



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 10:16 AM EDT

PDB ID : 4CY6 / pdb_00004cy6
Title : apo structure of 2-hydroxybiphenyl 3-monooxygenase HbpA
Authors : Jensen, C.N.; Farrugia, J.E.; Frank, A.; Man, H.; Hart, S.; Turkenburg, J.P.; Grogan, G.
Deposited on : 2014-04-10
Resolution : 2.76 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

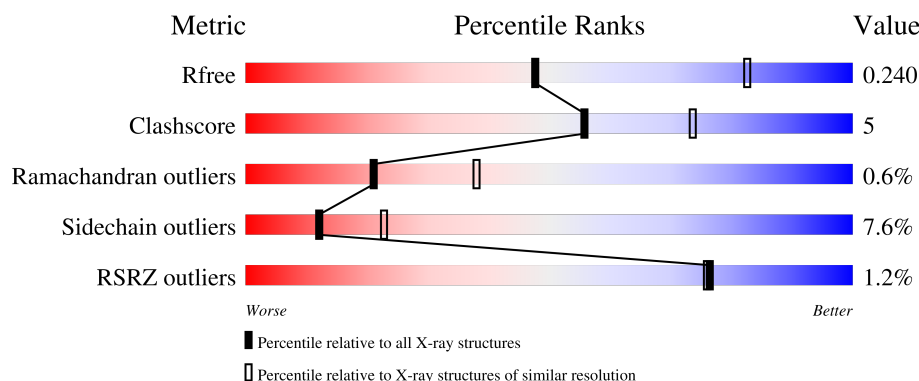
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	586	% 80% 13% • •
1	B	586	2% 80% 12% • 5%
1	C	586	2% 78% 13% • 5%
1	D	586	% 79% 13% • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16265 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 2-HYDROXYBIPHENYL-3-MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	561	Total	C	N	O	S	0	0	0
			4095	2603	723	749	20			
1	B	555	Total	C	N	O	S	0	0	0
			3982	2532	707	724	19			
1	C	554	Total	C	N	O	S	0	0	0
			4017	2545	717	736	19			
1	D	553	Total	C	N	O	S	0	0	0
			4032	2556	716	740	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	347	GLN	THR	engineered mutation	UNP O06647
B	347	GLN	THR	engineered mutation	UNP O06647
C	347	GLN	THR	engineered mutation	UNP O06647
D	347	GLN	THR	engineered mutation	UNP O06647

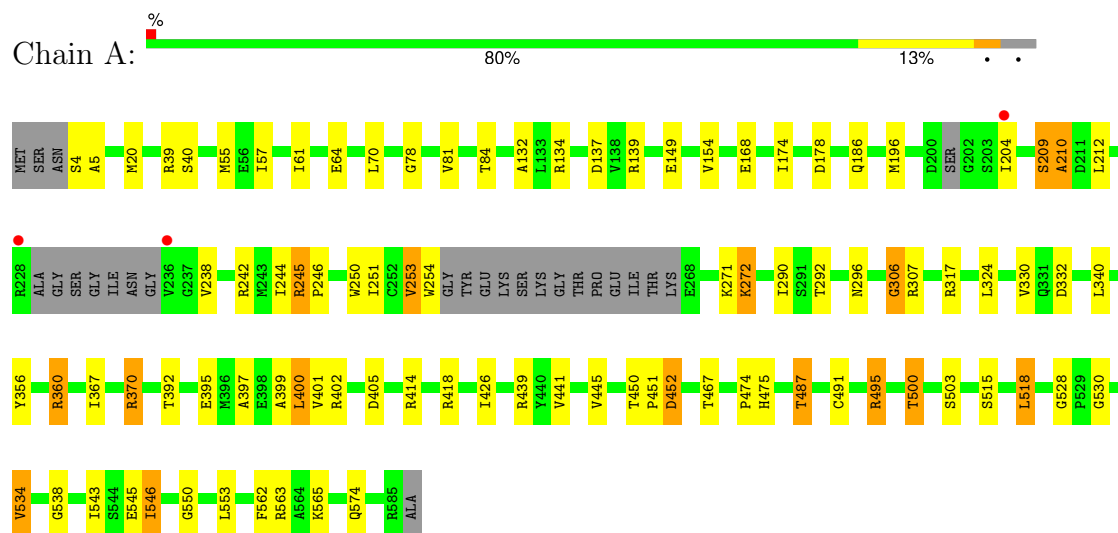
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	40	Total	O	0	0
			40	40		
2	B	35	Total	O	0	0
			35	35		
2	C	33	Total	O	0	0
			33	33		
2	D	31	Total	O	0	0
			31	31		

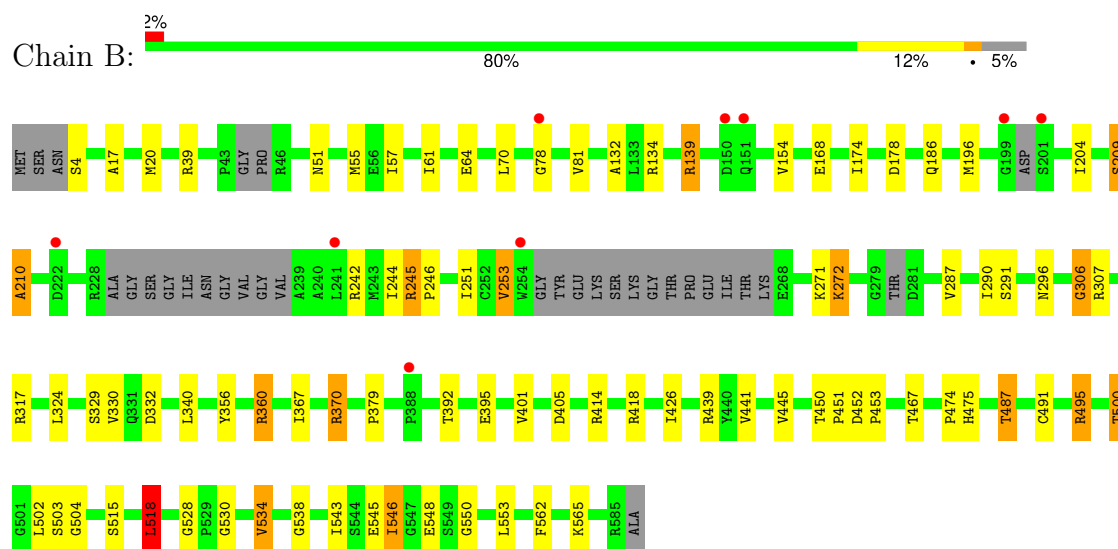
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

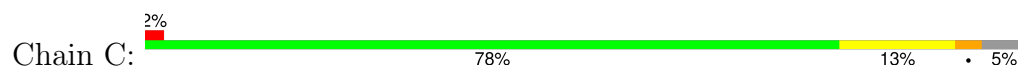
• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE

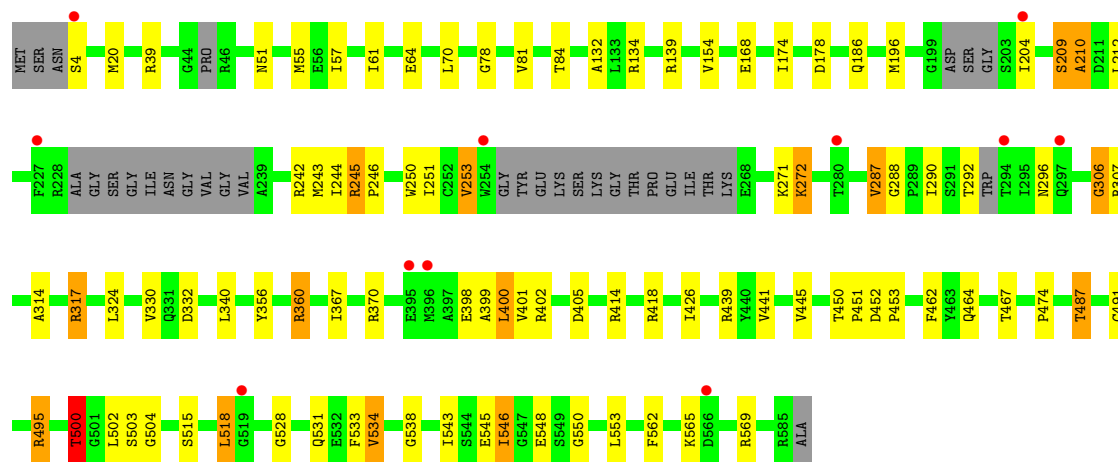


• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE

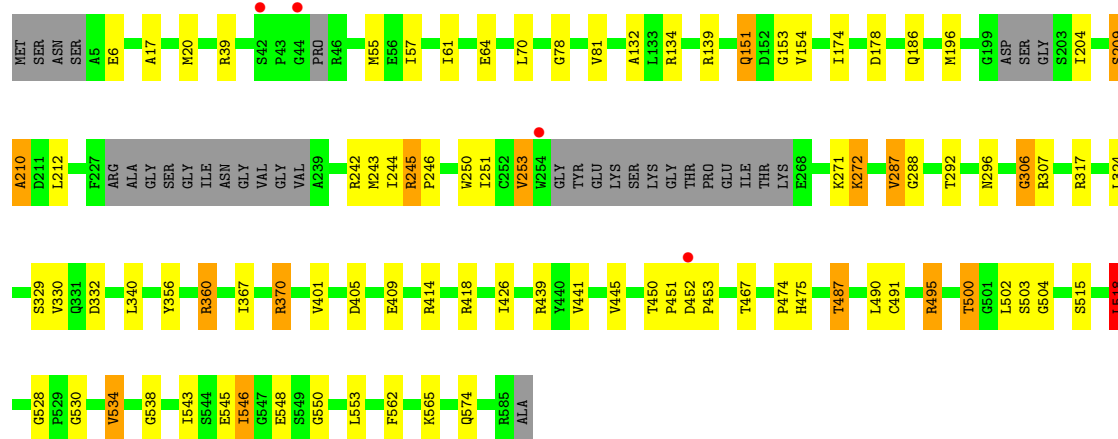
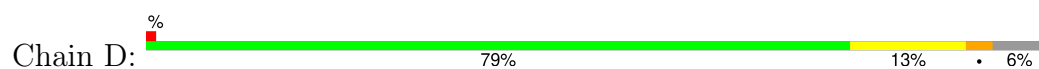


• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE





• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.35Å 94.67Å 99.16Å 115.03° 96.09° 109.59°	Depositor
Resolution (Å)	85.97 – 2.76 85.97 – 2.76	Depositor EDS
% Data completeness (in resolution range)	96.8 (85.97-2.76) 96.8 (85.97-2.76)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.190 , 0.241 0.195 , 0.240	Depositor DCC
R_{free} test set	2997 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.419	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16265	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.96	0/4184	1.06	13/5695 (0.2%)
1	B	0.98	0/4068	1.08	15/5547 (0.3%)
1	C	0.95	0/4102	1.07	15/5587 (0.3%)
1	D	0.97	2/4119 (0.0%)	1.06	14/5610 (0.2%)
All	All	0.97	2/16473 (0.0%)	1.07	57/22439 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	4
1	C	0	4
1	D	0	5
All	All	0	17

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	452	ASP	C-N	5.49	1.40	1.33
1	D	453	PRO	N-CD	5.20	1.55	1.47

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	253	VAL	N-CA-C	8.08	118.45	111.90
1	B	78	GLY	N-CA-C	7.92	122.91	112.77
1	C	528	GLY	CA-C-N	7.84	127.39	119.24
1	C	528	GLY	C-N-CA	7.84	127.39	119.24
1	B	253	VAL	N-CA-C	7.75	118.17	111.90
1	A	528	GLY	CA-C-N	7.55	127.09	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	528	GLY	C-N-CA	7.55	127.09	119.24
1	A	491	CYS	N-CA-C	7.29	124.96	114.39
1	C	491	CYS	N-CA-C	7.20	124.83	114.39
1	A	452	ASP	CA-C-N	-6.98	113.05	120.66
1	A	452	ASP	C-N-CA	-6.98	113.05	120.66
1	B	452	ASP	CA-C-N	-6.97	113.48	120.31
1	B	452	ASP	C-N-CA	-6.97	113.48	120.31
1	A	253	VAL	N-CA-C	6.91	117.50	111.90
1	B	453	PRO	N-CA-C	-6.84	101.18	111.57
1	D	253	VAL	N-CA-C	6.84	117.44	111.90
1	C	452	ASP	CA-C-N	-6.64	113.80	120.31
1	C	452	ASP	C-N-CA	-6.64	113.80	120.31
1	D	78	GLY	N-CA-C	6.64	121.27	112.77
1	B	491	CYS	N-CA-C	6.60	123.95	114.39
1	A	78	GLY	N-CA-C	6.57	121.18	112.77
1	A	238	VAL	N-CA-C	6.54	117.90	108.48
1	C	78	GLY	N-CA-C	6.44	121.01	112.77
1	D	528	GLY	CA-C-N	6.28	127.69	119.84
1	D	528	GLY	C-N-CA	6.28	127.69	119.84
1	D	452	ASP	CA-C-N	-6.25	112.64	120.51
1	D	452	ASP	C-N-CA	-6.25	112.64	120.51
1	B	528	GLY	CA-C-N	6.18	127.56	119.84
1	B	528	GLY	C-N-CA	6.18	127.56	119.84
1	D	491	CYS	N-CA-C	6.17	123.34	114.39
1	C	287	VAL	CB-CA-C	-5.82	104.27	111.08
1	D	534	VAL	CA-C-N	-5.68	114.40	122.41
1	D	534	VAL	C-N-CA	-5.68	114.40	122.41
1	C	518	LEU	N-CA-C	-5.58	101.77	110.42
1	D	409	GLU	N-CA-C	-5.50	105.37	111.36
1	C	534	VAL	CA-C-N	-5.48	114.65	122.77
1	C	534	VAL	C-N-CA	-5.48	114.65	122.77
1	A	534	VAL	CA-C-N	-5.47	114.67	122.77
1	A	534	VAL	C-N-CA	-5.47	114.67	122.77
1	D	287	VAL	CB-CA-C	-5.46	104.06	111.15
1	B	272	LYS	N-CA-C	-5.45	105.32	112.41
1	B	379	PRO	CA-C-N	-5.45	114.00	119.56
1	B	379	PRO	C-N-CA	-5.45	114.00	119.56
1	D	518	LEU	N-CA-C	-5.44	101.99	110.42
1	C	272	LYS	N-CA-C	-5.37	105.43	112.41
1	D	272	LYS	N-CA-C	-5.32	105.50	112.41
1	A	518	LEU	N-CA-C	-5.22	102.32	110.42
1	A	272	LYS	N-CA-C	-5.20	105.66	112.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	534	VAL	CA-C-N	-5.17	115.11	122.41
1	B	534	VAL	C-N-CA	-5.17	115.11	122.41
1	C	462	PHE	N-CA-C	5.13	117.26	108.90
1	C	453	PRO	N-CA-C	-5.08	103.84	111.57
1	B	518	LEU	N-CA-C	-5.07	102.56	110.42
1	A	40	SER	N-CA-C	5.06	115.72	108.74
1	B	287	VAL	CB-CA-C	-5.04	105.19	111.08
1	C	500	THR	CB-CA-C	-5.02	101.33	111.17
1	D	153	GLY	N-CA-C	5.00	116.95	110.45

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	271	LYS	Peptide
1	A	272	LYS	Peptide
1	A	530	GLY	Peptide
1	A	545	GLU	Peptide
1	B	271	LYS	Peptide
1	B	272	LYS	Peptide
1	B	530	GLY	Peptide
1	B	545	GLU	Peptide
1	C	271	LYS	Peptide
1	C	272	LYS	Peptide
1	C	531	GLN	Peptide
1	C	545	GLU	Peptide
1	D	271	LYS	Peptide
1	D	272	LYS	Peptide
1	D	490	LEU	Peptide
1	D	530	GLY	Peptide
1	D	545	GLU	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4095	0	3918	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3982	0	3750	40	0
1	C	4017	0	3798	42	0
1	D	4032	0	3830	38	0
2	A	40	0	0	2	0
2	B	35	0	0	0	0
2	C	33	0	0	2	0
2	D	31	0	0	0	0
All	All	16265	0	15296	149	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (149) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:ALA:O	1:B:39:ARG:NH1	1.97	0.96
1:A:196:MET:O	1:A:296:ASN:ND2	2.07	0.86
1:D:196:MET:O	1:D:296:ASN:ND2	2.08	0.86
1:B:196:MET:O	1:B:296:ASN:ND2	2.08	0.85
1:C:196:MET:O	1:C:296:ASN:ND2	2.11	0.83
1:B:405:ASP:O	1:B:414:ARG:NH2	2.13	0.82
1:C:39:ARG:NH1	1:D:132:ALA:O	2.14	0.81
1:A:405:ASP:O	1:A:414:ARG:NH2	2.13	0.80
1:A:39:ARG:NH1	1:B:132:ALA:O	2.14	0.80
1:D:405:ASP:O	1:D:414:ARG:NH2	2.14	0.80
1:D:70:LEU:O	1:D:245:ARG:NH2	2.16	0.79
1:C:405:ASP:O	1:C:414:ARG:NH2	2.16	0.79
1:C:70:LEU:O	1:C:245:ARG:NH2	2.18	0.77
1:B:70:LEU:O	1:B:245:ARG:NH2	2.17	0.76
1:B:515:SER:O	1:B:518:LEU:O	2.03	0.76
1:A:70:LEU:O	1:A:245:ARG:NH2	2.19	0.75
1:C:132:ALA:O	1:D:39:ARG:NH1	2.20	0.75
1:C:515:SER:O	1:C:518:LEU:O	2.06	0.74
1:A:515:SER:O	1:A:518:LEU:O	2.06	0.73
1:B:154:VAL:HG21	1:B:174:ILE:HG13	1.71	0.73
1:D:515:SER:O	1:D:518:LEU:O	2.06	0.72
1:D:154:VAL:HG21	1:D:174:ILE:HG13	1.74	0.69
1:C:154:VAL:HG21	1:C:174:ILE:HG13	1.74	0.69
1:D:244:ILE:O	1:D:245:ARG:HD2	1.93	0.69
1:A:154:VAL:HG21	1:A:174:ILE:HG13	1.75	0.68
1:C:244:ILE:O	1:C:245:ARG:HD2	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ILE:O	1:A:245:ARG:HD2	1.93	0.68
1:B:244:ILE:O	1:B:245:ARG:HD2	1.93	0.68
1:A:500:THR:HG23	1:A:550:GLY:O	1.96	0.65
1:B:500:THR:HG23	1:B:550:GLY:O	1.96	0.65
1:A:149:GLU:OE1	1:A:306:GLY:O	2.15	0.65
1:C:500:THR:HG23	1:C:550:GLY:O	1.98	0.64
1:D:306:GLY:O	1:D:307:ARG:HB2	1.98	0.64
1:A:306:GLY:O	1:A:307:ARG:HB2	1.97	0.63
1:B:306:GLY:O	1:B:307:ARG:HB2	1.99	0.63
1:D:500:THR:HG23	1:D:550:GLY:O	1.98	0.63
2:A:2016:HOH:O	1:B:139:ARG:HD3	1.98	0.62
1:B:414:ARG:NH1	1:D:548:GLU:OE2	2.33	0.62
1:C:306:GLY:O	1:C:307:ARG:HB2	1.98	0.61
1:A:84:THR:HG21	1:A:400:LEU:HD11	1.83	0.60
1:B:418:ARG:HD3	1:D:538:GLY:HA3	1.83	0.60
1:D:474:PRO:O	1:D:487:THR:HG21	2.02	0.59
1:B:474:PRO:O	1:B:487:THR:HG21	2.01	0.58
1:C:84:THR:HG21	1:C:400:LEU:HD11	1.85	0.58
1:A:474:PRO:O	1:A:487:THR:HG21	2.04	0.57
1:B:495:ARG:CG	1:B:495:ARG:HH11	2.18	0.57
1:A:495:ARG:CG	1:A:495:ARG:HH11	2.18	0.57
1:D:332:ASP:OD1	1:D:360:ARG:NH1	2.38	0.57
1:B:546:ILE:HD12	1:B:565:LYS:HA	1.87	0.56
1:C:474:PRO:O	1:C:487:THR:HG21	2.05	0.56
1:B:55:MET:HE3	1:B:64:GLU:HG3	1.88	0.55
1:C:495:ARG:HH11	1:C:495:ARG:CG	2.19	0.55
1:A:55:MET:HE3	1:A:64:GLU:HG3	1.88	0.55
1:C:55:MET:HE3	1:C:64:GLU:HG3	1.89	0.55
1:A:495:ARG:HH11	1:A:495:ARG:HG3	1.72	0.54
1:B:332:ASP:OD1	1:B:360:ARG:NH1	2.41	0.54
1:D:55:MET:HE3	1:D:64:GLU:HG3	1.89	0.54
1:C:332:ASP:OD1	1:C:360:ARG:NH1	2.41	0.54
1:C:546:ILE:HD12	1:C:565:LYS:HA	1.90	0.53
1:A:332:ASP:OD1	1:A:360:ARG:NH1	2.39	0.53
1:A:414:ARG:NH1	1:C:548:GLU:OE2	2.42	0.53
1:D:495:ARG:HH11	1:D:495:ARG:HG3	1.74	0.53
1:C:495:ARG:HH11	1:C:495:ARG:HG3	1.73	0.53
1:B:495:ARG:HH11	1:B:495:ARG:HG3	1.73	0.52
1:D:495:ARG:HH11	1:D:495:ARG:CG	2.22	0.52
1:D:450:THR:HG22	1:D:451:PRO:O	2.11	0.51
1:B:548:GLU:OE2	1:D:414:ARG:NH1	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:546:ILE:HD12	1:D:565:LYS:HA	1.94	0.50
1:A:553:LEU:HB3	1:A:562:PHE:HB3	1.93	0.49
1:A:5:ALA:N	2:A:2001:HOH:O	2.46	0.49
1:A:204:ILE:O	1:A:253:VAL:O	2.31	0.48
1:A:546:ILE:HD12	1:A:565:LYS:HA	1.95	0.48
1:C:553:LEU:HB3	1:C:562:PHE:HB3	1.96	0.48
1:D:204:ILE:O	1:D:253:VAL:O	2.32	0.47
1:B:553:LEU:HB3	1:B:562:PHE:HB3	1.96	0.47
1:A:137:ASP:OD1	1:B:39:ARG:HD3	2.15	0.47
1:A:209:SER:O	1:A:210:ALA:HB2	2.15	0.47
1:D:360:ARG:HH11	1:D:360:ARG:HG3	1.80	0.47
1:D:553:LEU:HB3	1:D:562:PHE:HB3	1.96	0.47
1:C:209:SER:O	1:C:210:ALA:HB2	2.15	0.47
1:D:151:GLN:HE21	1:D:151:GLN:HA	1.80	0.47
1:B:204:ILE:O	1:B:253:VAL:O	2.33	0.46
1:B:209:SER:O	1:B:210:ALA:HB2	2.15	0.46
1:B:475:HIS:HA	1:B:487:THR:HG22	1.98	0.46
1:C:399:ALA:HA	1:C:402:ARG:HD3	1.96	0.46
1:A:399:ALA:HA	1:A:402:ARG:HD3	1.98	0.45
1:B:20:MET:HG2	1:B:330:VAL:HG13	1.97	0.45
1:C:51:ASN:OD1	1:C:51:ASN:C	2.60	0.45
1:D:209:SER:O	1:D:210:ALA:HB2	2.16	0.45
1:D:20:MET:HG2	1:D:330:VAL:HG13	1.97	0.45
1:B:154:VAL:HG21	1:B:174:ILE:CG1	2.45	0.45
1:C:204:ILE:O	1:C:253:VAL:O	2.34	0.45
1:D:370:ARG:NH2	1:D:426:ILE:HG12	2.32	0.45
1:A:20:MET:HG2	1:A:330:VAL:HG13	1.98	0.44
1:B:242:ARG:HD3	1:B:251:ILE:HD11	2.00	0.44
1:B:370:ARG:NH2	1:B:426:ILE:HG12	2.32	0.44
1:C:243:MET:C	1:C:243:MET:SD	3.01	0.44
1:C:20:MET:HG2	1:C:330:VAL:HG13	2.00	0.44
1:B:502:LEU:C	1:B:504:GLY:H	2.26	0.44
1:A:450:THR:HG22	1:A:451:PRO:O	2.17	0.43
1:C:398:GLU:O	1:C:402:ARG:HD3	2.19	0.43
1:B:450:THR:HG22	1:B:451:PRO:O	2.19	0.43
1:B:392:THR:HG22	1:B:395:GLU:OE1	2.19	0.43
1:C:450:THR:HG22	1:C:451:PRO:O	2.18	0.43
1:C:502:LEU:C	1:C:504:GLY:H	2.26	0.43
1:D:212:LEU:HD12	1:D:250:TRP:CE2	2.53	0.43
1:A:57:ILE:O	1:A:61:ILE:HG12	2.19	0.43
1:D:242:ARG:HD3	1:D:251:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:ARG:HD3	1:C:538:GLY:HA3	1.99	0.42
1:C:287:VAL:HG12	1:C:288:GLY:O	2.20	0.42
1:A:538:GLY:HA3	1:C:418:ARG:HD3	2.00	0.42
1:C:360:ARG:NH2	2:C:2021:HOH:O	2.41	0.42
1:A:356:TYR:O	1:A:360:ARG:HB2	2.20	0.42
1:C:370:ARG:NH2	1:C:426:ILE:HG12	2.35	0.42
1:A:475:HIS:HA	1:A:487:THR:HG22	2.01	0.42
1:C:4:SER:HA	1:C:168:GLU:O	2.20	0.42
1:A:212:LEU:HD12	1:A:250:TRP:CE2	2.55	0.42
1:B:495:ARG:CG	1:B:495:ARG:NH1	2.83	0.42
1:C:242:ARG:HD3	1:C:251:ILE:HD11	2.00	0.42
1:D:57:ILE:O	1:D:61:ILE:HG12	2.19	0.42
1:B:4:SER:HA	1:B:168:GLU:O	2.20	0.42
1:C:464:GLN:HG2	2:C:2031:HOH:O	2.19	0.42
1:A:242:ARG:HD3	1:A:251:ILE:HD11	2.02	0.42
1:A:360:ARG:HH11	1:A:360:ARG:HG3	1.85	0.42
1:B:61:ILE:HG22	1:B:61:ILE:O	2.20	0.42
1:D:61:ILE:O	1:D:61:ILE:HG22	2.20	0.42
1:A:392:THR:HG22	1:A:395:GLU:OE1	2.20	0.41
1:D:475:HIS:HA	1:D:487:THR:HG22	2.02	0.41
1:D:502:LEU:C	1:D:504:GLY:H	2.29	0.41
1:A:546:ILE:CG1	1:A:563:ARG:HD2	2.50	0.41
1:C:314:ALA:O	1:C:317:ARG:NH1	2.54	0.41
1:A:370:ARG:NH2	1:A:426:ILE:HG12	2.36	0.41
1:B:51:ASN:OD1	1:B:51:ASN:C	2.64	0.41
1:A:4:SER:HA	1:A:168:GLU:O	2.21	0.41
1:A:154:VAL:HG21	1:A:174:ILE:CG1	2.49	0.41
1:B:17:ALA:HB2	1:B:329:SER:HB3	2.03	0.41
1:B:57:ILE:O	1:B:61:ILE:HG12	2.21	0.41
1:B:538:GLY:HA3	1:D:418:ARG:HD3	2.02	0.41
1:C:356:TYR:O	1:C:360:ARG:HB2	2.21	0.41
1:C:360:ARG:HH11	1:C:360:ARG:HG3	1.85	0.41
1:D:17:ALA:HB2	1:D:329:SER:HB3	2.03	0.41
1:D:287:VAL:HG12	1:D:288:GLY:O	2.21	0.41
1:C:57:ILE:O	1:C:61:ILE:HG12	2.21	0.40
1:C:212:LEU:HD12	1:C:250:TRP:CE2	2.55	0.40
1:D:243:MET:SD	1:D:243:MET:C	3.04	0.40
1:D:356:TYR:O	1:D:360:ARG:HB2	2.21	0.40
1:A:397:ALA:O	1:A:401:VAL:HG13	2.21	0.40
1:B:356:TYR:O	1:B:360:ARG:HB2	2.21	0.40
1:C:154:VAL:HG21	1:C:174:ILE:CG1	2.47	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	553/586 (94%)	521 (94%)	29 (5%)	3 (0%)	24	43
1	B	543/586 (93%)	514 (95%)	26 (5%)	3 (1%)	21	38
1	C	542/586 (92%)	510 (94%)	28 (5%)	4 (1%)	18	34
1	D	543/586 (93%)	513 (94%)	27 (5%)	3 (1%)	21	38
All	All	2181/2344 (93%)	2058 (94%)	110 (5%)	13 (1%)	21	38

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	GLY
1	B	306	GLY
1	C	306	GLY
1	D	306	GLY
1	A	210	ALA
1	B	210	ALA
1	C	210	ALA
1	C	533	PHE
1	D	210	ALA
1	A	246	PRO
1	B	246	PRO
1	C	246	PRO
1	D	246	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	392/467 (84%)	362 (92%)	30 (8%)	12	22
1	B	371/467 (79%)	343 (92%)	28 (8%)	12	24
1	C	379/467 (81%)	351 (93%)	28 (7%)	13	24
1	D	385/467 (82%)	355 (92%)	30 (8%)	11	21
All	All	1527/1868 (82%)	1411 (92%)	116 (8%)	12	23

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	VAL
1	A	134	ARG
1	A	139	ARG
1	A	178	ASP
1	A	186	GLN
1	A	209	SER
1	A	245	ARG
1	A	254	TRP
1	A	290	ILE
1	A	292	THR
1	A	317	ARG
1	A	324	LEU
1	A	340	LEU
1	A	360	ARG
1	A	367	ILE
1	A	370	ARG
1	A	400	LEU
1	A	439	ARG
1	A	441	VAL
1	A	445	VAL
1	A	452	ASP
1	A	467	THR
1	A	487	THR
1	A	495	ARG
1	A	500	THR
1	A	503	SER
1	A	534	VAL
1	A	543	ILE
1	A	546	ILE
1	A	574	GLN

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Mol	Chain	Res	Type
1	B	81	VAL
1	B	134	ARG
1	B	139	ARG
1	B	178	ASP
1	B	186	GLN
1	B	209	SER
1	B	245	ARG
1	B	290	ILE
1	B	291	SER
1	B	317	ARG
1	B	324	LEU
1	B	340	LEU
1	B	360	ARG
1	B	367	ILE
1	B	370	ARG
1	B	401	VAL
1	B	439	ARG
1	B	441	VAL
1	B	445	VAL
1	B	467	THR
1	B	487	THR
1	B	495	ARG
1	B	500	THR
1	B	503	SER
1	B	518	LEU
1	B	534	VAL
1	B	543	ILE
1	B	546	ILE
1	C	81	VAL
1	C	134	ARG
1	C	139	ARG
1	C	178	ASP
1	C	186	GLN
1	C	209	SER
1	C	245	ARG
1	C	290	ILE
1	C	292	THR
1	C	317	ARG
1	C	324	LEU
1	C	340	LEU
1	C	360	ARG
1	C	367	ILE

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Mol	Chain	Res	Type
1	C	400	LEU
1	C	401	VAL
1	C	439	ARG
1	C	441	VAL
1	C	445	VAL
1	C	467	THR
1	C	487	THR
1	C	495	ARG
1	C	500	THR
1	C	503	SER
1	C	534	VAL
1	C	543	ILE
1	C	546	ILE
1	C	569	ARG
1	D	6	GLU
1	D	81	VAL
1	D	134	ARG
1	D	139	ARG
1	D	151	GLN
1	D	178	ASP
1	D	186	GLN
1	D	209	SER
1	D	245	ARG
1	D	292	THR
1	D	317	ARG
1	D	324	LEU
1	D	340	LEU
1	D	360	ARG
1	D	367	ILE
1	D	370	ARG
1	D	401	VAL
1	D	439	ARG
1	D	441	VAL
1	D	445	VAL
1	D	467	THR
1	D	487	THR
1	D	495	ARG
1	D	500	THR
1	D	503	SER
1	D	518	LEU
1	D	534	VAL
1	D	543	ILE

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Mol	Chain	Res	Type
1	D	546	ILE
1	D	574	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	HIS
1	A	275	HIS
1	A	459	GLN
1	B	107	HIS
1	C	107	HIS
1	C	275	HIS
1	C	481	ASN
1	C	516	GLN
1	D	100	HIS
1	D	107	HIS
1	D	151	GLN
1	D	275	HIS
1	D	459	GLN
1	D	481	ASN
1	D	509	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	561/586 (95%)	-0.14	3 (0%) 87 87	39, 59, 81, 118	0
1	B	555/586 (94%)	-0.07	9 (1%) 70 69	39, 61, 86, 119	0
1	C	554/586 (94%)	-0.01	11 (1%) 65 63	40, 60, 87, 113	0
1	D	553/586 (94%)	-0.13	4 (0%) 84 84	37, 60, 86, 117	0
All	All	2223/2344 (94%)	-0.09	27 (1%) 76 76	37, 60, 86, 119	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	VAL	4.6
1	B	222	ASP	3.5
1	C	294	THR	3.5
1	D	254	TRP	3.4
1	C	227	PHE	3.2
1	B	254	TRP	3.2
1	A	228	ARG	3.1
1	D	44	GLY	2.9
1	C	4	SER	2.9
1	C	395	GLU	2.8
1	C	204	ILE	2.7
1	D	42	SER	2.7
1	B	151	GLN	2.6
1	B	241	LEU	2.6
1	C	280	THR	2.5
1	C	396	MET	2.5
1	B	201	SER	2.3
1	C	519	GLY	2.3
1	B	150	ASP	2.3
1	A	204	ILE	2.2
1	C	254	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	452	ASP	2.2
1	B	388	PRO	2.1
1	C	297	GLN	2.1
1	B	78	GLY	2.1
1	B	199	GLY	2.1
1	C	566	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.